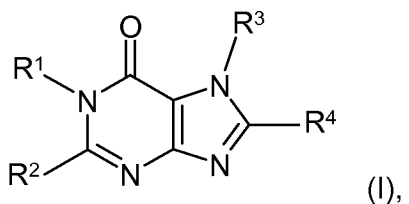


AMENDMENTS TO THE CLAIMS

1. (previously presented) A compound of the formula (I):



wherein

R¹ denotes a hydrogen atom,

a C₁₋₈-alkyl group,

a C₃₋₈-alkenyl group,

a C₃₋₄-alkenyl group which is substituted by a C₁₋₂-alkyloxy-carbonyl, aminocarbonyl, C₁₋₃-alkylamino-carbonyl, di-(C₁₋₃-alkyl)-amino-carbonyl, pyrrolidin-1-ylcarbonyl, piperidin-1-ylcarbonyl or morpholin-4-ylcarbonyl group,

a C₃₋₈-alkynyl group,

a C₁₋₆-alkyl group substituted by a group R_a, where

R_a denotes a C₃₋₇-cycloalkyl, heteroaryl, cyano, carboxy, C₁₋₃-alkyloxy-carbonyl, aminocarbonyl, C₁₋₃-alkylamino-carbonyl, di-(C₁₋₃-alkyl)-amino-carbonyl, pyrrolidin-1-ylcarbonyl, piperidin-1-ylcarbonyl, morpholin-4-ylcarbonyl, piperazin-1-ylcarbonyl, 4-methylpiperazin-1-ylcarbonyl or 4-ethylpiperazin-1-ylcarbonyl group,

a C₁₋₆-alkyl group substituted by a phenyl group, where the phenyl ring is substituted by the groups R¹⁰ to R¹⁴ and

R¹⁰ denotes a hydrogen atom,

a fluorine, chlorine, bromine or iodine atom,

a C₁₋₄-alkyl, hydroxy or C₁₋₄-alkyloxy group,

a nitro, amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, cyan-C₁₋₃-alkylamino, N-(cyan-C₁₋₃-alkyl)-N-(C₁₋₃-alkyl)-amino, C₁₋₃-alkyloxy-carbonyl-C₁₋₃-alkylamino, pyrrolidin-1-yl, piperidin-1-yl, morpholin-4-yl, piperazin-1-yl or 4-(C₁₋₃-alkyl)-piperazin-1-yl group,

a formylamino, C₁₋₃-alkyl-carbonylamino, C₃₋₆-cycloalkyl-carbonylamino, C₃₋₆-cycloalkyl-C₁₋₃-alkyl-carbonylamino, arylcarbonylamino, aryl-C₁₋₃-alkyl-carbonylamino, C₁₋₃-alkyloxy-carbonylamino, aminocarbonylamino, C₁₋₃-alkyl-aminocarbonylamino, di-(C₁₋₃-alkyl)-aminocarbonylamino, pyrrolidin-1-yl-carbonylamino, piperidin-1-yl-carbonylamino, morpholin-4-yl-carbonylamino,

piperazin-1-yl-carbonylamino or 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonylamino, C₁₋₃-alkyl-sulphonylamino, bis-(C₁₋₃-alkylsulphonyl)-amino, aminosulphonylamino, C₁₋₃-alkylamino-sulphonylamino, di-(C₁₋₃-alkyl)-amino-sulphonylamino, pyrrolidin-1-yl-sulphonylamino, piperidin-1-yl-sulphonylamino, morpholin-4-yl-sulphonylamino, piperazin-1-yl-sulphonylamino or 4-(C₁₋₃-alkyl)-piperazin-1-yl-sulphonylamino, (C₁₋₃-alkylamino)-thiocarbonylamino, (C₁₋₃-alkyloxy-carbonylamino)-carbonylamino, arylsulphonylamino or aryl-C₁₋₃-alkyl-sulphonylamino group,

an N-(C₁₋₃-alkyl)-formylamino, N-(C₁₋₃-alkyl)-N-(C₁₋₃-alkyl-carbonyl)-amino, N-(C₁₋₃-alkyl)-N-(C₃₋₆-cycloalkyl-carbonyl)-amino, N-(C₁₋₃-alkyl)-N-(C₃₋₆-cycloalkyl-C₁₋₃-alkyl-carbonyl)-amino, N-(C₁₋₃-alkyl)-N-(arylcarbonyl)-amino, N-(C₁₋₃-alkyl)-N-(aryl-C₁₋₃-alkyl-carbonyl)-amino, N-(C₁₋₃-alkyl)-N-(C₁₋₃-alkyloxy-carbonyl)-amino, N-(aminocarbonyl)-N-(C₁₋₃-alkyl)-amino, N-(C₁₋₃-alkyl-amino-carbonyl)-N-(C₁₋₃-alkyl)-amino, N-[di-(C₁₋₃-alkyl)aminocarbonyl]-N-(C₁₋₃-alkyl)-amino, N-(C₁₋₃-alkyl)-N-(C₁₋₃-alkyl-sulphonyl)-amino, N-(C₁₋₃-alkyl)-N-(aryl-sulphonyl)-amino or N-(C₁₋₃-alkyl)-N-(aryl-C₁₋₃-alkyl-sulphonyl)-amino group,

a 2-oxo-imidazolidin-1-yl, 2,4-dioxo-imidazolidin-1-yl, 2,5-dioxo-imidazolidin-1-yl or 2-oxo-hexahydropyrimidin-1-yl group wherein the nitrogen atom in the 3 position may be substituted in each case by a methyl or ethyl group,

a cyano, carboxy, C₁₋₃-alkyloxy-carbonyl, aminocarbonyl, C₁₋₃-alkyl-amino-carbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, pyrrolidin-1-yl-carbonyl, piperidin-1-yl-

carbonyl, morpholin-4-yl-carbonyl, piperazin-1-yl-carbonyl or 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl group,

a C₁₋₃-alkyl-carbonyl or an arylcarbonyl group,

a carboxy-C₁₋₃-alkyl, C₁₋₃-alkyloxy-carbonyl-C₁₋₃-alkyl, cyano-C₁₋₃-alkyl, aminocarbonyl-C₁₋₃-alkyl, C₁₋₃-alkyl-aminocarbonyl-C₁₋₃-alkyl, di-(C₁₋₃-alkyl)-aminocarbonyl-C₁₋₃-alkyl, pyrrolidin-1-yl-carbonyl-C₁₋₃-alkyl, piperidin-1-yl-carbonyl-C₁₋₃-alkyl, morpholin-4-yl-carbonyl-C₁₋₃-alkyl, piperazin-1-yl-carbonyl-C₁₋₃-alkyl or 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl-C₁₋₃-alkyl group,

a carboxy-C₁₋₃-alkyloxy, C₁₋₃-alkyloxy-carbonyl-C₁₋₃-alkyloxy, cyano-C₁₋₃-alkyloxy, aminocarbonyl-C₁₋₃-alkyloxy, C₁₋₃-alkyl-aminocarbonyl-C₁₋₃-alkyloxy, di-(C₁₋₃-alkyl)-aminocarbonyl-C₁₋₃-alkyloxy, pyrrolidin-1-yl-carbonyl-C₁₋₃-alkyloxy, piperidin-1-yl-carbonyl-C₁₋₃-alkyloxy, morpholin-4-yl-carbonyl-C₁₋₃-alkyloxy, piperazin-1-yl-carbonyl-C₁₋₃-alkyloxy or 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl-C₁₋₃-alkyloxy group,

a hydroxy-C₁₋₃-alkyl, C₁₋₃-alkyloxy-C₁₋₃-alkyl, amino-C₁₋₃-alkyl, C₁₋₃-alkylamino-C₁₋₃-alkyl, di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl, pyrrolidin-1-yl-C₁₋₃-alkyl, piperidin-1-yl-C₁₋₃-alkyl, morpholin-4-yl-C₁₋₃-alkyl, piperazin-1-yl-C₁₋₃-alkyl, 4-(C₁₋₃-alkyl)-piperazin-1-yl-C₁₋₃-alkyl group,

a hydroxy-C₁₋₃-alkyloxy, C₁₋₃-alkyloxy-C₁₋₃-alkyloxy, C₁₋₃-alkylsulphanyl-C₁₋₃-alkyloxy, C₁₋₃-alkylsulphinyl-C₁₋₃-alkyloxy, C₁₋₃-alkylsulphonyl-C₁₋₃-alkyloxy,

amino- C_{1-3} -alkyloxy, C_{1-3} -alkylamino- C_{1-3} -alkyloxy, di-(C_{1-3} -alkyl)-amino- C_{1-3} -alkyloxy, pyrrolidin-1-yl- C_{1-3} -alkyloxy, piperidin-1-yl- C_{1-3} -alkyloxy, morpholin-4-yl- C_{1-3} -alkyloxy, piperazin-1-yl- C_{1-3} -alkyloxy, 4-(C_{1-3} -alkyl)-piperazin-1-yl- C_{1-3} -alkyloxy group,

a mercapto, C_{1-3} -alkylsulphanyl, C_{1-3} -alkylsulphinyl, C_{1-3} -alkylsulphonyl, C_{1-3} -alkylsulphonyloxy, arylsulphonyloxy, trifluoromethylsulphanyl, trifluoromethylsulphinyl or trifluoromethylsulphonyl group,

a sulpho, aminosulphonyl, C_{1-3} -alkyl-aminosulphonyl, di-(C_{1-3} -alkyl)-aminosulphonyl, pyrrolidin-1-yl-sulphonyl, piperidin-1-yl-sulphonyl, morpholin-4-yl-sulphonyl, piperazin-1-yl-sulphonyl or 4-(C_{1-3} -alkyl)-piperazin-1-yl-sulphonyl group,

a methyl or methoxy group substituted by 1 to 3 fluorine atoms,

an ethyl or ethyloxy group substituted by 1 to 5 fluorine atoms,

a C_{2-4} -alkenyl or C_{2-4} -alkynyl group,

a C_{3-4} -alkenyloxy or C_{3-4} -alkynyloxy group,

a C_{3-6} -cycloalkyl or C_{3-6} -cycloalkyloxy group,

a C_{3-6} -cycloalkyl- C_{1-3} -alkyl or C_{3-6} -cycloalkyl- C_{1-3} -alkyloxy group or

an aryl, aryloxy, aryl-C₁₋₃-alkyl or aryl-C₁₋₃-alkyloxy group,

R¹¹ and R¹², which may be identical or different, in each case represent a hydrogen atom, a fluorine, chlorine, bromine or iodine atom, a C₁₋₃-alkyl, trifluoromethyl, hydroxy, C₁₋₃-alkyloxy or cyano group, or

R¹¹ together with R¹², if they are bound to adjacent carbon atoms, also represent a methylenedioxy, difluoromethylenedioxy or a straight-chain C₃₋₅-alkylene group and

R¹³ and R¹⁴, which may be identical or different, in each case represent a hydrogen atom, a fluorine, chlorine or bromine atom, a trifluoromethyl, C₁₋₃-alkyl or C₁₋₃-alkyloxy group,

a phenyl-C₁₋₄-alkyl group wherein the alkyl moiety is substituted by a cyano, carboxy, C₁₋₃-alkyloxy-carbonyl, aminocarbonyl, C₁₋₃-alkyl-aminocarbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, pyrrolidin-1-yl-carbonyl, piperidin-1-yl-carbonyl, morpholin-4-yl-carbonyl-group and the phenyl moiety is substituted by the groups R¹⁰ to R¹⁴, while R¹⁰ to R¹⁴ are as hereinbefore defined,

a phenyl group substituted by the groups R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴ are as hereinbefore defined,

a phenyl-C₂₋₃-alkenyl group wherein the phenyl moiety is substituted by the groups R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴ are as hereinbefore defined,

a phenyl-(CH₂)_m-A-(CH₂)_n-group wherein the phenyl moiety is substituted by R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴ are as hereinbefore defined and

A represents a carbonyl group, m represents the number 0, 1 or 2 and n represents the number 1, 2 or 3,

a phenylcarbonylmethyl group wherein the phenyl moiety is substituted by R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴ are as hereinbefore defined and the methyl moiety is substituted by a C₁₋₃-alkyl group,

a phenyl-(CH₂)_m-B-(CH₂)_n-group wherein the phenyl moiety is substituted by R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴, m and n are as hereinbefore defined and

B denotes a methylene group which is substituted by a hydroxy, C₁₋₃-alkyloxy, amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, mercapto, C₁₋₃-alkylsulphanyl, C₁₋₃-alkylsulphinyl or C₁₋₃-alkylsulphonyl group and is optionally additionally substituted by a methyl or ethyl group,

a naphthyl-C₁₋₃-alkyl group wherein the naphthyl moiety is substituted by the groups R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴ are as hereinbefore defined,

a naphthyl-(CH₂)_m-A-(CH₂)_n-group wherein the naphthyl moiety is substituted by R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴, A, m and n are as hereinbefore defined,

a naphthyl-(CH₂)_m-B-(CH₂)_n-group wherein the naphthyl moiety is substituted by R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴, B, m and n are as hereinbefore defined,

a [1,4]naphthoquinon-2-yl, chromen-4-on-3-yl, 1-oxoindan-2-yl, 1,3-dioxoindan-2-yl or 2,3-dihydro-3-oxo-benzofuran-2-yl group,

a heteroaryl-(CH₂)_m-A-(CH₂)_n group where A, m and n are as hereinbefore defined,

a heteroaryl-(CH₂)_m-B-(CH₂)_n group where B, m and n are as hereinbefore defined,

a C₁₋₆-alkyl-A-(CH₂)_n group where A and n are as hereinbefore defined,

a C₃₋₇-cycloalkyl-(CH₂)_m-A-(CH₂)_n group where A, m and n are as hereinbefore defined,

a C₃₋₇-cycloalkyl-(CH₂)_m-B-(CH₂)_n group where B, m and n are as hereinbefore defined,

an R²¹-A-(CH₂)_n-group wherein R²¹ denotes a C₁₋₃-alkyloxycarbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl, di-(C₁₋₃-alkyl)aminocarbonyl, pyrrolidin-1-yl-carbonyl, piperidin-1-yl-carbonyl or morpholin-4-yl-carbonyl, piperazin-1-yl-carbonyl, 4-

methylpiperazin-1-yl-carbonyl or 4-ethylpiperazin-1-yl-carbonyl group and A and n are as hereinbefore defined,

a phenyl-(CH₂)_m-D-C₁₋₃-alkyl group wherein the phenyl moiety is substituted by the groups R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴ and m are as mentioned hereinbefore and D denotes an oxygen or sulphur atom, -NH-, C₁₋₃-alkylimino, sulphinyl or sulphonyl group,

a naphthyl-(CH₂)_m-D-C₁₋₃-alkyl group wherein the naphthyl moiety is substituted by the groups R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴, D and m are as mentioned hereinbefore,

a C₂₋₆-alkyl group substituted by a group R_b, where

R_b is isolated from the cyclic nitrogen atom in the 1 position of the purine skeleton by at least two carbon atoms and

R_b denotes a hydroxy, C₁₋₃-alkyloxy, mercapto, C₁₋₃-alkylsulphanyl, C₁₋₃-alkylsulphinyl, C₁₋₃-alkylsulphonyl, amino, C₁₋₃-alkyl-carbonylamino, C₃₋₆-cycloalkyl-carbonyl-amino, arylcarbonylamino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, pyrrolidin-1-yl, piperidin-1-yl, morpholin-4-yl, piperazin-1-yl or 4-(C₁₋₃-alkyl)-piperazin-1-yl group,

a C₃₋₆-cycloalkyl group,

or an amino or arylcarbonylamino group,

R^2 denotes a hydrogen atom,

a C_{1-8} -alkyl group,

a C_{3-8} -alkenyl group,

a C_{3-4} -alkenyl group which is substituted by a C_{1-2} -alkyloxy-carbonyl, aminocarbonyl, C_{1-3} -alkylamino-carbonyl, di- $(C_{1-3}$ -alkyl)-amino-carbonyl, pyrrolidin-1-ylcarbonyl, piperidin-1-ylcarbonyl or morpholin-4-ylcarbonyl group,

a C_{3-8} -alkynyl group,

a C_{3-6} -cycloalkyl group,

a C_{1-6} -alkyl group substituted by a group R_a , where R_a is as hereinbefore defined,

a phenyl group which is substituted by R^{10} to R^{14} , where R^{10} to R^{14} are as hereinbefore defined,

a C_{1-6} -alkyl group substituted by a phenyl group, wherein the phenyl ring is substituted by the groups R^{10} to R^{14} and R^{10} to R^{14} are as hereinbefore defined,

a phenyl- C_{1-4} -alkyl group wherein the alkyl moiety is substituted by a cyano, carboxy, C_{1-3} -alkyloxy-carbonyl, aminocarbonyl, C_{1-3} -alkyl-aminocarbonyl, di- $(C_{1-3}$ -alkyl)-

aminocarbonyl, pyrrolidin-1-yl-carbonyl, piperidin-1-yl-carbonyl, morpholin-4-yl-carbonyl group and the phenyl moiety is substituted by the groups R^{10} to R^{14} , where R^{10} to R^{14} are as hereinbefore defined,

a phenyl- C_{2-3} -alkenyl group wherein the phenyl moiety is substituted by R^{10} to R^{14} , where R^{10} to R^{14} are as hereinbefore defined,

a heteroaryl group,

a phenyl- $(CH_2)_m$ -A or phenyl- $(CH_2)_m$ -A- $(CH_2)_n$ group wherein the phenyl moiety is substituted in each case by R^{10} to R^{14} , while A, R^{10} to R^{14} , m and n are as hereinbefore defined,

a phenylcarbonylmethyl group wherein the phenyl moiety is substituted by R^{10} to R^{14} , where R^{10} to R^{14} are as hereinbefore defined and the methyl moiety is substituted by a C_{1-3} -alkyl group,

a phenyl- $(CH_2)_m$ -B or phenyl- $(CH_2)_m$ -B- $(CH_2)_n$ group wherein the phenyl moiety is substituted in each case by R^{10} to R^{14} , while B, R^{10} to R^{14} , m and n are as hereinbefore defined,

a naphthyl- C_{1-3} -alkyl group wherein the naphthyl moiety is substituted by the groups R^{10} to R^{14} , where R^{10} to R^{14} are as hereinbefore defined,

a naphthyl-(CH₂)_m-A or naphthyl-(CH₂)_m-A-(CH₂)_n group wherein the naphthyl moiety is substituted in each case by R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴, A, m and n are as hereinbefore defined,

a naphthyl-(CH₂)_m-B or naphthyl-(CH₂)_m-B-(CH₂)_n group wherein the naphthyl moiety is substituted in each case by R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴, B, m and n are as hereinbefore defined,

a heteroaryl-(CH₂)_m-A or heteroaryl-(CH₂)_m-A-(CH₂)_n group where A, m and n are as hereinbefore defined,

a heteroaryl-(CH₂)_m-B or heteroaryl-(CH₂)_m-B-(CH₂)_n group where B, m and n are as hereinbefore defined,

a C₁₋₆-alkyl-A or C₁₋₆-alkyl-A-(CH₂)_n group where A and n are as hereinbefore defined,

a C₃₋₇-cycloalkyl-(CH₂)_m-A or C₃₋₇-cycloalkyl-(CH₂)_m-A-(CH₂)_n group where A, m and n are as hereinbefore defined,

a C₃₋₇-cycloalkyl-(CH₂)_m-B or C₃₋₇-cycloalkyl-(CH₂)_m-B-(CH₂)_n group where B, m and n are as hereinbefore defined,

an R²¹-A-(CH₂)_n group wherein R²¹, A and n are as hereinbefore defined,

a phenyl-(CH₂)_m-D-C₁₋₃-alkyl group wherein the phenyl moiety is substituted by the groups R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴, D and m are as mentioned hereinbefore,

a naphthyl-(CH₂)_m-D-C₁₋₃-alkyl group wherein the naphthyl moiety is substituted by the groups R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴, D and m are as mentioned hereinbefore,

a C₁₋₆-alkyl group substituted by a group R_b, where R_b is as hereinbefore defined,

a cyano, carboxy, C₁₋₃-alkyloxy-carbonyl, aminocarbonyl, C₁₋₃-alkylamino-carbonyl, di-(C₁₋₃-alkyl)-amino-carbonyl, pyrrolidin-1-ylcarbonyl, piperidin-1-ylcarbonyl, morpholin-4-ylcarbonyl, piperazin-1-ylcarbonyl, 4-methylpiperazin-1-ylcarbonyl or 4-ethylpiperazin-1-ylcarbonyl group,

an amino, C₁₋₆-alkylamino or di-(C₁₋₆-alkyl)-amino group,

an amino group substituted by the groups R¹⁵ and R¹⁶ wherein

R¹⁵ denotes a hydrogen atom or a C₁₋₆-alkyl group and

R¹⁶ denotes a C₁₋₆-alkyl group which is substituted by R_a, where R_a is as hereinbefore defined,

an amino group substituted by the groups R¹⁵ and R¹⁷ wherein

R¹⁵ is as hereinbefore defined and

R¹⁷ denotes a C₂₋₆-alkyl group which is substituted by a hydroxy, C₁₋₃-alkyloxy, aryloxy, mercapto, C₁₋₃-alkylsulphanyl, C₁₋₃-alkylsulphinyl, C₁₋₃-alkylsulphonyl, C₁₋₃-alkylsulphonylamino, arylsulphanyl, arylsulphinyl, arylsulphonyl, arylsulphonylamino, C₁₋₃-alkyl-carbonylamino, C₃₋₆-cycloalkyl-carbonylamino, arylcarbonylamino, C₁₋₃-alkyl-oxycarbonylamino, aminocarbonylamino, C₁₋₃-alkyl-aminocarbonylamino, di-(C₁₋₃-alkyl)-aminocarbonylamino, amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, pyrrolidin-1-yl, piperidin-1-yl, morpholin-4-yl, piperazin-1-yl or 4-(C₁₋₃-alkyl)-piperazin-1-yl group,

a C₃₋₆-cycloalkylamino or N-(C₃₋₆-cycloalkyl)-N-(C₁₋₃-alkyl)-amino group,

a phenylamino or N-(phenyl)-N-(C₁₋₃-alkyl)-amino group wherein the phenyl moiety is substituted in each case by R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴ are as hereinbefore defined,

a phenyl-C₁₋₆-alkylamino or N-(phenyl-C₁₋₆-alkyl)-N-(C₁₋₃-alkyl)-amino group wherein the phenyl moiety is substituted in each case by R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴ are as hereinbefore defined,

a naphthylamino or N-(naphthyl)-N-(C₁₋₃-alkyl)-amino group,

a naphthyl-C₁₋₆-alkylamino or N-(naphthyl-C₁₋₆-alkyl)-N-(C₁₋₃-alkyl)-amino group,

a heteroaryl amino or N-(heteroaryl)-N-(C₁₋₃-alkyl)-amino group,

a pyrrolidin-1-yl, piperidin-1-yl, homopiperidin-1-yl, morpholin-4-yl, homomorpholin-4-yl, piperazin-1-yl, 4-(C₁₋₃-alkyl)-piperazin-1-yl, homopiperazin-1-yl or 4-(C₁₋₃-alkyl)-homopiperazin-1-yl group, or

a C₁₋₆-alkyloxy, C₃₋₆-cycloalkyloxy or C₃₋₆-cycloalkyl-C₁₋₆-alkyloxy group,

a C₁₋₆-alkylsulphanyl, C₃₋₆-cycloalkylsulphanyl or C₃₋₆-cycloalkyl-C₁₋₆-alkylsulphanyl group,

a phenyloxy or phenylsulphanyl group wherein the phenyl moiety is substituted in each case by R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴ are as hereinbefore defined,

a phenyl-C₁₋₆-alkyloxy or phenyl-C₁₋₆-alkylsulphanyl group wherein the phenyl moiety is substituted in each case by R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴ are as hereinbefore defined,

a naphthyloxy or a naphthylsulphanyl group wherein the naphthyl moiety is substituted in each case by R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴ are as hereinbefore defined,

a naphthyl-C₁₋₆-alkyloxy or naphthyl-C₁₋₆-alkylsulphanyl group wherein the naphthyl moiety is substituted in each case by R¹⁰ to R¹⁴, where R¹⁰ to R¹⁴ are as hereinbefore defined,

a heteroaryloxy or heteroarylsulphanyl group or

a heteroaryl-C₁₋₆-alkyloxy or heteroaryl-C₁₋₆-alkylsulphanyl group,

R³ denotes a C₁₋₈-alkyl group,

a C₁₋₄-alkyl group substituted by the group R_c, where

R_c denotes a C₃₋₇-cycloalkyl group optionally substituted by one or two C₁₋₃-alkyl groups,

a C₅₋₇-cycloalkenyl group optionally substituted by one or two C₁₋₃-alkyl groups,

an aryl group or

a furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, pyridyl, pyridazinyl, pyrimidyl or pyrazinyl group, while the above-mentioned heterocyclic groups may each be substituted by one or two C₁₋₃-alkyl groups or by a fluorine, chlorine, bromine or iodine atom or by a trifluoromethyl, cyano or C₁₋₃-alkyloxy group,

a C₃₋₈-alkenyl group,

a C₃₋₆-alkenyl group substituted by a fluorine, chlorine or bromine atom or a trifluoromethyl group,

a C₃₋₈-alkynyl group,

an aryl group or

an aryl-C₂₋₄-alkenyl group,

and

R⁴ denotes an azetidin-1-yl or pyrrolidin-1-yl group which is substituted in the 3 position by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group and may additionally be substituted by one or two C₁₋₃-alkyl groups,

a piperidin-1-yl or hexahydroazepin-1-yl group which is substituted in the 3 position or in the 4 position by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)amino group and may additionally be substituted by one or two C₁₋₃-alkyl groups,

a 3-amino-piperidin-1-yl group wherein the piperidin-1-yl-moiety is additionally substituted by an aminocarbonyl, C₁₋₂-alkyl-aminocarbonyl, di-(C₁₋₂-alkyl)-aminocarbonyl, pyrrolidin-1-yl-carbonyl, (2-cyano-pyrrolidin-1-yl)-carbonyl, thiazolidin-3-yl-carbonyl, (4-cyano-thiazolidin-3-yl)carbonyl, piperidin-1-ylcarbonyl or morpholin-4-ylcarbonyl group,

a 3-amino-piperidin-1-yl group wherein the piperidin-1-yl moiety is additionally substituted in the 4 position or 5 position by a hydroxy or methoxy group,

a 3-amino-piperidin-1-yl group wherein the methylene group is replaced in the 2 position or 6 position by a carbonyl group,

a piperidin-1-yl or hexahydroazepin-1-yl group substituted in the 3 position by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group, wherein two hydrogen atoms on the carbon skeleton of the piperidin-1-yl or hexahydroazepin-1-yl group are each replaced by a straight-chain alkylene bridge, this bridge containing 2 to 5 carbon atoms if the two hydrogen atoms are on the same carbon atom, or 1 to 4 carbon atoms if the hydrogen atoms are on adjacent carbon atoms, or 1 to 4 carbon atoms if the hydrogen atoms are on carbon atoms which are separated by one atom, or 1 to 3 carbon atoms if the hydrogen atoms are on carbon atoms which are separated by two atoms,

an azetidin-1-yl, pyrrolidin-1-yl, piperidin-1-yl or hexahydroazepin-1-yl group which is substituted by an amino-C₁₋₃-alkyl, C₁₋₃-alkylamino-C₁₋₃-alkyl or a di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl group,

a piperazin-1-yl or [1,4]diazepan-1-yl group optionally substituted on the carbon skeleton by one or two C₁₋₃-alkyl groups,

a 3-imino-piperazin-1-yl, 3-imino-[1,4]diazepan-1-yl or 5-imino-[1,4]diazepan-1-yl group optionally substituted on the carbon skeleton by one or two C₁₋₃-alkyl groups,

a [1,4]diazepan-1-yl group optionally substituted by one or two C₁₋₃-alkyl groups, which is substituted in the 6 position by an amino group,

a C₃₋₇-cycloalkyl group which is substituted by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group,

a C₃₋₇-cycloalkyl group which is substituted by an amino-C₁₋₃-alkyl, C₁₋₃-alkylamino-C₁₋₃-alkyl or a di-(C₁₋₃-alkyl)amino-C₁₋₃-alkyl group,

a C₃₋₇-cycloalkyl-C₁₋₂-alkyl group wherein the cycloalkyl moiety is substituted by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group,

a C₃₋₇-cycloalkyl-C₁₋₂-alkyl group wherein the cycloalkyl moiety is substituted by an amino-C₁₋₃-alkyl, C₁₋₃-alkylamino-C₁₋₃-alkyl or a di-(C₁₋₃-alkyl)amino-C₁₋₃-alkyl group,

a C₃₋₇-cycloalkylamino group wherein the cycloalkyl moiety is substituted by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group, while the two nitrogen atoms on the cycloalkyl moiety are separated from one another by at least two carbon atoms,

an N-(C₃₋₇-cycloalkyl)-N-(C₁₋₃-alkyl)-amino group wherein the cycloalkyl moiety is substituted by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group, while the two nitrogen atoms on the cycloalkyl moiety are separated from one another by at least two carbon atoms,

a C₃₋₇-cycloalkylamino group wherein the cycloalkyl moiety is substituted by an amino-C₁₋₃-alkyl, C₁₋₃-alkylamino-C₁₋₃-alkyl or a di-(C₁₋₃-alkyl)amino-C₁₋₃-alkyl group,

an N-(C₃₋₇-cycloalkyl)-N-(C₁₋₃-alkyl)-amino group wherein the cycloalkyl moiety is substituted by an amino-C₁₋₃-alkyl, C₁₋₃-alkylamino-C₁₋₃-alkyl or a di-(C₁₋₃-alkyl)amino-C₁₋₃-alkyl group,

a C₃₋₇-cycloalkyl-C₁₋₂-alkyl-amino group wherein the cycloalkyl moiety is substituted by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group,

an N-(C₃₋₇-cycloalkyl-C₁₋₂-alkyl)-N-(C₁₋₂-alkyl)-amino group wherein the cycloalkyl moiety is substituted by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group,

a C₃₋₇-cycloalkyl-C₁₋₂-alkyl-amino group wherein the cycloalkyl moiety is substituted by an amino-C₁₋₃-alkyl, C₁₋₃-alkylamino-C₁₋₃-alkyl or a di-(C₁₋₃-alkyl)amino-C₁₋₃-alkyl group,

an N-(C₃₋₇-cycloalkyl-C₁₋₂-alkyl)-N-(C₁₋₂-alkyl)-amino group wherein the cycloalkyl moiety is substituted by an amino-C₁₋₃-alkyl, C₁₋₃-alkylamino-C₁₋₃-alkyl or a di-(C₁₋₃-alkyl)amino-C₁₋₃-alkyl group,

an R¹⁹-C₂₋₄-alkylamino group wherein R¹⁹ is separated from the nitrogen atom of the C₂₋₄-alkylamino moiety by at least two carbon atoms and

R¹⁹ denotes an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group,

an R¹⁹-C₂₋₄-alkylamino group wherein the nitrogen atom of the C₂₋₄-alkylamino moiety is substituted by a C₁₋₃-alkyl group and R¹⁹ is separated from the nitrogen atom of the C₂₋₄-alkylamino moiety by at least two carbon atoms, where R¹⁹ is as hereinbefore defined,

an amino group substituted by the group R²⁰ wherein

R²⁰ denotes an azetidin-3-yl, azetidin-2-ylmethyl, azetidin-3-ylmethyl, pyrrolidin-3-yl, pyrrolidin-2-ylmethyl, pyrrolidin-3-ylmethyl, piperidin-3-yl, piperidin-4-yl, piperidin-2-ylmethyl, piperidin-3-ylmethyl or piperidin-4-ylmethyl group, while the groups mentioned for R²⁰ may each be substituted by one or two C₁₋₃-alkyl groups,

an amino group substituted by the group R²⁰ and a C₁₋₃-alkyl group wherein R²⁰ is as hereinbefore defined, while the groups mentioned for R²⁰ may each be substituted by one or two C₁₋₃-alkyl groups,

an R¹⁹-C₃₋₄-alkyl group wherein the C₃₋₄-alkyl moiety is straight-chained and may additionally be substituted by one or two C₁₋₃-alkyl groups, where R¹⁹ is as hereinbefore defined,

a 3-amino-2-oxo-piperidin-5-yl or 3-amino-2-oxo-1-methyl-piperidin-5-yl group,

a pyrrolidin-3-yl, piperidin-3-yl, piperidin-4-yl, hexahydroazepin-3-yl or hexahydroazepin-4-yl group which is substituted in the 1 position by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)amino group,

or an azetidin-2-yl-C₁₋₂-alkyl, azetidin-3-yl-C₁₋₂-alkyl, pyrrolidin-2-yl-C₁₋₂-alkyl, pyrrolidin-3-yl, pyrrolidin-3-yl-C₁₋₂-alkyl, piperidin-2-yl-C₁₋₂-alkyl, piperidin-3-yl, piperidin-3-yl-C₁₋₂-alkyl, piperidin-4-yl or piperidin-4-yl-C₁₋₂-alkyl group, while the above-mentioned groups may each be substituted by one or two C₁₋₃-alkyl groups,

while by the aryl groups mentioned in the definition of the above groups are meant phenyl or naphthyl groups, which may be mono- or disubstituted by R_h independently of one another, where the substituents are identical or different and R_h denotes a fluorine, chlorine, bromine or iodine atom, a trifluoromethyl, cyano, nitro, amino, aminocarbonyl, aminosulphonyl, methylsulphonyl, acetylamino, methylsulphonylamino, C₁₋₃-alkyl, cyclopropyl, ethenyl, ethynyl, hydroxy, C₁₋₃-alkyloxy, difluoromethoxy or trifluoromethoxy group,

by the heteroaryl groups mentioned in the definitions of the above-mentioned groups are meant a pyrrolyl, furanyl, thienyl, pyridyl, indolyl, benzofuranyl, benzothiophenyl, quinolinyl or isoquinolinyl group,

or a pyrrolyl, furanyl, thienyl or pyridyl group wherein one or two methyne groups are replaced by nitrogen atoms,

or an indolyl, benzofuranyl, benzothiophenyl, quinolinyl or isoquinolinyl group wherein one to three methyne groups are replaced by nitrogen atoms,

or a 1,2-dihydro-2-oxo-pyridinyl, 1,4-dihydro-4-oxo-pyridinyl, 2,3-dihydro-3-oxo-pyridazinyl, 1,2,3,6-tetrahydro-3,6-dioxo-pyridazinyl, 1,2-dihydro-2-oxo-pyrimidinyl, 3,4-dihydro-4-oxo-pyrimidinyl, 1,2,3,4-tetrahydro-2,4-dioxo-pyrimidinyl, 1,2-dihydro-2-oxo-pyrazinyl, 1,2,3,4-tetrahydro-2,3-dioxo-pyrazinyl, 2,3-dihydro-2-oxo-indolyl, 2,3-dihydrobenzofuranyl, 2,3-dihydro-2-oxo-1H-benzimidazolyl, 2,3-dihydro-2-oxo-benzoxazolyl, 1,2-dihydro-2-oxo-quinolinyl, 1,4-dihydro-4-oxo-quinolinyl, 1,2-dihydro-1-oxo-isoquinolinyl, 1,4-dihydro-4-oxo-cinnolinyl, 1,2-dihydro-2-oxo-quinazolinyl, 3,4-dihydro-4-oxo-quinazolinyl, 1,2,3,4-tetrahydro-2,4-dioxo-quinazolinyl, 1,2-dihydro-2-oxoquinoxalinyl, 1,2,3,4-tetrahydro-2,3-dioxo-quinoxalinyl, 1,2-dihydro-1-oxo-phthalazinyl, 1,2,3,4-tetrahydro-1,4-dioxo-phthalazinyl, chromanyl, cumaranyl, 2,3-dihydro-benzo[1,4]dioxinyl or 3,4-dihydro-3-oxo-2H-benzo[1,4]oxazinyl group,

while the above-mentioned heteroaryl groups may be substituted by R^{10} to R^{14} , where R^{10} to R^{14} are as hereinbefore defined,

and, unless otherwise stated, the above-mentioned alkyl, alkenyl and alkynyl groups may be straight-chain or branched,

as well as the derivatives which are N-oxidised at the cyclic nitrogen atom in the 3 position or 9 position of the hypoxanthine skeleton,

as well as the derivatives wherein the 6-oxo group of the hypoxanthine skeleton is replaced by a thioxo group,

with the proviso that the compounds

8-(piperidin-4-ylmethyl)-7-(4-fluorobenzyl)-1,7-dihydro-purin-6-one and

8-(1-methyl-piperidin-4-ylmethyl)-7-(4-fluorobenzyl)-1,7-dihydro-purin-6-one

are excluded,

the tautomers, enantiomers, diastereomers, the mixtures thereof, the prodrugs thereof and the salts thereof.

2. (original) The compound according to claim 1,

wherein R^1 , R^2 and R^3 are defined as in claim 1 and

R^4 denotes a pyrrolidin-1-yl group which is substituted in the 3 position by an amino group,

a piperidin-1-yl group which is substituted in the 3 position by an amino group,

a piperidin-3-yl or piperidin-4-yl group,

a hexahydroazepin-1-yl group which is substituted in the 3 position or 4 position by an amino group,

a piperazin-1-yl or [1,4]diazepan-1-yl group,

a (2-aminocyclohexyl)amino group,

a cyclohexyl group which is substituted in the 3 position by an amino group,

or an N-(2-aminoethyl)-methylamino or an N-(2-aminoethyl)-ethylamino group,

the tautomers, enantiomers, diastereomers, the mixtures thereof, the prodrugs and the salts thereof.

3. (original) The compound according to claim 1, wherein

R¹ denotes a hydrogen atom,

a C₁₋₆-alkyl group,

a C₃₋₆-alkenyl group,

a C₃₋₄-alkynyl group,

a C₃₋₆-cycloalkylmethyl group,

a phenyl-C₁₋₃-alkyl group wherein the phenyl moiety is substituted by R¹⁰ and R¹¹,
where

R¹⁰ denotes a hydrogen atom, a fluorine, chlorine or bromine atom,

a methyl or trifluoromethyl group,

a cyano, aminocarbonyl, dimethylaminocarbonyl or methylsulphonyl group,

an amino, acetylamino or methylsulphonylamino group,

a hydroxy, methoxy, difluoromethoxy, trifluoromethoxy, carboxymethoxy,
methoxycarbonylmethoxy, ethyloxycarbonylmethoxy, aminocarbonylmethoxy,
methylaminocarbonylmethoxy, ethylaminocarbonylmethoxy or
dimethylaminocarbonylmethoxy group and

R¹¹ denotes a hydrogen atom, a fluorine or chlorine atom,

or a methyl or methoxy group,

a naphthylmethyl group wherein the naphthyl moiety is substituted by R¹⁰ and R¹¹,
where R¹⁰ and R¹¹ are as hereinbefore defined,

a heteroarylmethyl group where the term

heteroaryl denotes a furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, pyridyl, pyrimidinyl, pyrazinyl, quinoliny, isoquinoliny or quinazolinyl group and the above-mentioned heteroaryl groups are substituted by R^{10} and R^{11} , where R^{10} and R^{11} are as hereinbefore defined,

a phenylcarbonylmethyl group wherein the phenyl moiety is substituted by R^{10} and R^{11} , where R^{10} and R^{11} are as hereinbefore defined,

a furanylcabonylmethyl, thienylcarbonylmethyl or pyridylcarbonylmethyl group,

or a 2-oxo-propyl or cyclohexylcarbonylmethyl group,

R^2 denotes a hydrogen atom,

a C_{1-6} -alkyl group,

a C_{3-6} -alkenyl group,

a C_{3-4} -alkynyl group,

a C_{3-6} -cycloalkyl or C_{3-6} -cycloalkyl- C_{1-3} -alkyl group,

a phenyl group which is substituted by R^{10} and R^{11} , where R^{10} and R^{11} are as hereinbefore defined,

a phenyl-C₁₋₃-alkyl group wherein the phenyl moiety is substituted by R¹⁰ and R¹¹,
where R¹⁰ and R¹¹ are as hereinbefore defined,

a phenyl-C₂₋₃-alkenyl group wherein the phenyl moiety is substituted by R¹⁰ and R¹¹,
where R¹⁰ and R¹¹ are as hereinbefore defined,

a phenylcarbonylmethyl group wherein the phenyl moiety is substituted by R¹⁰ and
R¹¹, where R¹⁰ and R¹¹ are as hereinbefore defined,

a furanyl, thienyl or pyridyl group,

a furanyl-C₁₋₃-alkyl, thienyl-C₁₋₃-alkyl or pyridyl-C₁₋₃-alkyl group,

a cyano group,

an amino, C₁₋₄-alkylamino or di-(C₁₋₄-alkyl)-amino group,

an amino group substituted by the groups R¹⁵ and R¹⁶ wherein

R¹⁵ denotes a hydrogen atom or a methyl or ethyl group and

R¹⁶ denotes a C₁₋₄-alkyl group which is substituted by a cyano, carboxy,
methoxycarbonyl, ethyloxycarbonyl, aminocarbonyl, methylaminocarbonyl,

dimethylaminocarbonyl, ethylaminocarbonyl, diethylaminocarbonyl, pyrrolidin-1-ylcarbonyl or morpholin-4-ylcarbonyl group,

an amino group substituted by the groups R^{15} and R^{17} wherein

R^{15} is as hereinbefore defined and

R^{17} denotes a straight-chain C_{2-4} -alkyl group which is terminally substituted in each case by an amino, methylamino, dimethylamino, acetylamino, ethyloxy-carbonylamino, phenylcarbonylamino, methylsulphonylamino, phenylsulphonylamino, hydroxy, methoxy, phenyloxy, methylsulphanyl or phenylsulphanyl group,

a pyrrolidin-1-yl, piperidin-1-yl, morpholin-4-yl, piperazin-1-yl or 4-methyl-piperazin-1-yl group,

a C_{3-6} -cycloalkylamino or C_{3-6} -cycloalkyl- C_{1-3} -alkylamino group,

a phenylamino group,

a phenyl- C_{1-3} -alkylamino group wherein the phenyl moiety is substituted by R^{10} and R^{11} , where R^{10} and R^{11} are as hereinbefore defined,

a naphthylmethylamino group,

a heteroaryl- C_{1-2} -alkylamino group, where the term heteroaryl is as hereinbefore defined, or

a methylsulphanyl, benzylsulphanyl or (2-phenylethyl)sulphanyl group,

R^3 denotes a C_{4-6} -alkenyl group,

a C_{3-4} -alkenyl group which is substituted by a fluorine, chlorine or bromine atom or a trifluoromethyl group,

a 2-butyne-1-yl group or

a methyl group substituted by the group R_c , where

R_c denotes a 1-cyclopenten-1-yl-or 1-cyclohexen-1-yl group,

a phenyl group optionally substituted by a fluorine, chlorine, bromine or iodine atom, by a methyl, trifluoromethyl, cyano, methoxy, difluoromethoxy or trifluoromethoxy group,

a phenyl group which is substituted by two fluorine atoms,

a naphthyl group or

a furanyl, thienyl, or pyridyl group,

and

R^4 denotes a piperidin-1-yl group which is substituted in the 3 position by an amino group,

a hexahydroazepin-1-yl group which is substituted in the 3 position or 4 position by an amino group,

a (2-aminocyclohexyl)amino group,

a cyclohexyl group which is substituted in the 3 position by an amino group, or

an N-(2-aminoethyl)-methylamino or an N-(2-aminoethyl)-ethylamino group,

while unless otherwise stated, the above-mentioned alkyl, alkenyl and alkynyl groups may be straight-chain or branched.

4. (original) The compound according to claim 3, wherein

R^1 denotes a hydrogen atom,

a methyl, benzyl or 2-phenylethyl group,

a naphthylmethyl or methoxynaphthylmethyl group or

a phenylcarbonylmethyl group,

R^2 denotes a hydrogen atom,

a methyl or 2-phenylethyl group,

a phenylcarbonylmethyl group,

a cyano group,

an amino, methylamino, dimethylamino, isopropylamino, cyclohexylamino-
or (cyclohexylmethyl)amino group,

a benzylamino, fluorobenzylamino or (2-phenylethyl)amino group or

a piperidin-1-yl group,

R^3 denotes a benzyl or 3-methyl-but-2-en-1-yl group

and

R^4 denotes a (3-amino-piperidin-1-yl) group.

5. (original) A compound chosen from:

- (1) 8-(3-amino-piperidin-1-yl)-7-benzyl-2-benzylamino-1-methyl-1,7-dihydro-purin-6-one,
- (2) 8-(3-amino-piperidin-1-yl)-7-benzyl-2-(4-fluoro-benzylamino)-1-methyl-1,7-dihydro-purin-6-one,
- (3) 8-(3-amino-piperidin-1-yl)-7-benzyl-1-methyl-2-[(2-phenylethyl)amino]-1,7-dihydro-purin-6-one,
- (4) 8-(3-amino-piperidin-1-yl)-7-benzyl-2-isopropylamino-1-methyl-1,7-dihydro-purin-6-one,
- (5) 8-(3-amino-piperidin-1-yl)-7-benzyl-1-methyl-2-methylamino-1,7-dihydro-purin-6-one,
- (6) 8-(3-amino-piperidin-1-yl)-7-benzyl-2-cyclohexylamino-1-methyl-1,7-dihydro-purin-6-one,
- (7) 8-(3-amino-piperidin-1-yl)-7-benzyl-2-[(cyclohexylmethyl)amino]-1-methyl-1,7-dihydro-purin-6-one,
- (8) 8-(3-amino-piperidin-1-yl)-7-benzyl-1-methyl-2-(piperidin-1-yl)-1,7-dihydro-purin-6-one,

- (9) 8-(3-amino-piperidin-1-yl)-7-benzyl-2-dimethylamino-1-methyl-1,7-dihydro-purin-6-one,
- (10) 2-amino-8-(3-amino-piperidin-1-yl)-7-benzyl-1-methyl-1,7-dihydro-purin-6-one,
- (11) 8-(3-amino-piperidin-1-yl)-2-benzylamino-1-methyl-7-(3-methyl-but-2-en-1-yl)-1,7-dihydro-purin-6-one,
- (12) 8-(3-amino-piperidin-1-yl)-7-benzyl-1-methyl-2-methyl-1,7-dihydro-purin-6-one,
- (13) 8-(3-amino-piperidin-1-yl)-1-methyl-7-(3-methyl-but-2-en-1-yl)-2-(2-phenylethyl)-1,7-dihydro-purin-6-one,
- (14) 8-(3-amino-piperidin-1-yl)-7-benzyl-1-methyl-1,7-dihydro-purin-6-one,
- (15) 8-(3-amino-piperidin-1-yl)-7-benzyl-1-(2-oxo-2-phenyl-ethyl)-1,7-dihydro-purin-6-one,
- (16) 8-(3-amino-piperidin-1-yl)-2-methyl-7-(3-methyl-but-2-en-1-yl)-1-[(naphthalen-1-yl)methyl]-1,7-dihydro-purin-6-one,
- (17) 8-(3-amino-piperidin-1-yl)-7-(3-methyl-but-2-en-1-yl)-1-[(naphthalen-1-yl)methyl]-1,7-dihydro-purin-6-one and

(18) 8-(3-amino-piperidin-1-yl)-7-(3-methyl-but-2-en-1-yl)-1-[(4-methoxynaphthalen-1-yl)methyl]-1,7-dihydro-purin-6-one

as well as the tautomers, enantiomers, diastereomers, the mixtures thereof and the salts thereof.

6. (original) A physiologically acceptable salt of a compound according to claim 1 with inorganic or organic acids or bases.

7. (original) A pharmaceutical composition comprising a pharmaceutically effective amount of a compound according to claim 1 optionally together with one or more pharmaceutically acceptable inert carriers and/or diluents.

8. (currently amended) ~~A method of treating type I and type II diabetes mellitus, rheumatoid arthritis, obesity, allograft transplantation osteoporosis, or calcitonin-induced osteoporosis, said method~~ comprising administering to a patient in need thereof ~~a pharmaceutically effective amount of a compound according to claim 1~~ in an amount effective for the prevention or treatment of a disease or a condition associated with an increased DPP-IV activity.

9. (new) A method comprising administering to a patient in need thereof a compound according to claim 1 in an amount effective for the prevention or treatment of a disease or a condition that is capable of being prevented or alleviated by reducing the DPP-IV activity.